

Data Path : Z:\HPCHEM1\BNA E\DATA\BE103117\
 Data File : BE094681.D
 Acq On : 1 Nov 2017 4:53
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.470

Quant Time: Nov 01 09:13:44 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 08:49:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1306	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7692	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	4170	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9093	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	9421	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	8899	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4865	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	11188	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	7304	0.388	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	5095	0.387	ng/ul	98
7) Acenaphthylene	14.24	152	7681	0.407	ng/ul	97
8) Acenaphthene	14.58	153	5407	0.399	ng/ul	97
9) Fluorene	15.57	166	6148	0.404	ng/ul	93
11) Pentachlorophenol	17.01	266	171	0.377	ng/ul#	84
12) Phenanthrene	17.29	178	9858	0.410	ng/ul#	86
13) Anthracene	17.39	178	9548	0.403	ng/ul#	93
15) Fluoranthene	19.31	202	11348	0.409	ng/ul	84
17) Pyrene	19.68	202	12104	0.416	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	11206	0.416	ng/ul#	87
19) Chrysene	21.46	228	10860	0.390	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	10305	0.412	ng/ul	90
22) Benzo(k)fluoranthene	23.23	252	10653	0.397	ng/ul	92
23) Benzo(a)pyrene	23.86	252	10146	0.405	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.67	276	9140	0.353	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.67	278	7985	0.366	ng/ul	93
26) Benzo(g,h,i)perylene	27.52	276	6702	0.322	ng/ul	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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